



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:25 PM UTC

PDB ID : 2O0O / pdb_00002o0o
Title : Crystal structure of TL1A
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Deposited on : 2006-11-27
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

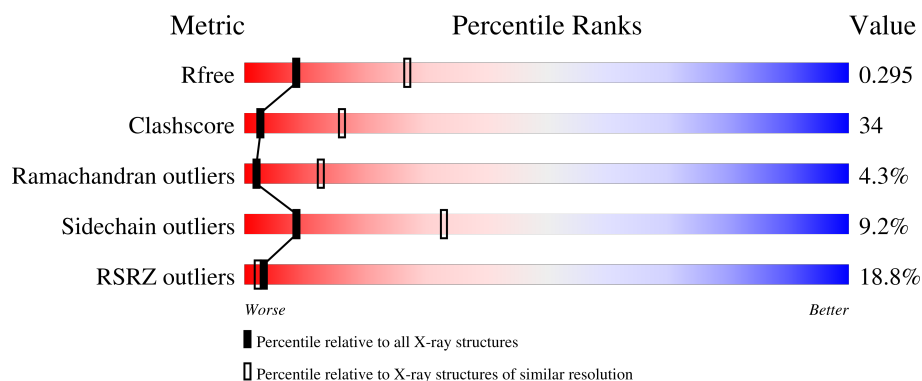
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	180	12% 36% 42% 6% 15%
1	B	180	15% 38% 38% 8% 15%
1	C	180	21% 33% 45% 6% 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3591 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called TNF superfamily ligand TL1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1200	777	194	222	7			
1	B	153	Total	C	N	O	S	0	0	0
			1183	764	190	223	6			
1	C	152	Total	C	N	O	S	0	0	0
			1170	757	188	219	6			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

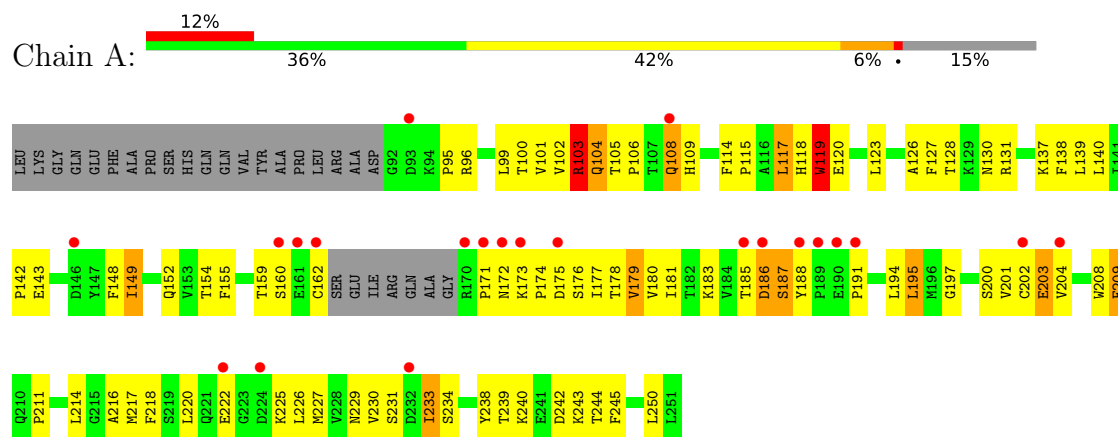
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	16	Total	O	0	0
			16	16		
3	B	12	Total	O	0	0
			12	12		
3	C	9	Total	O	0	0
			9	9		

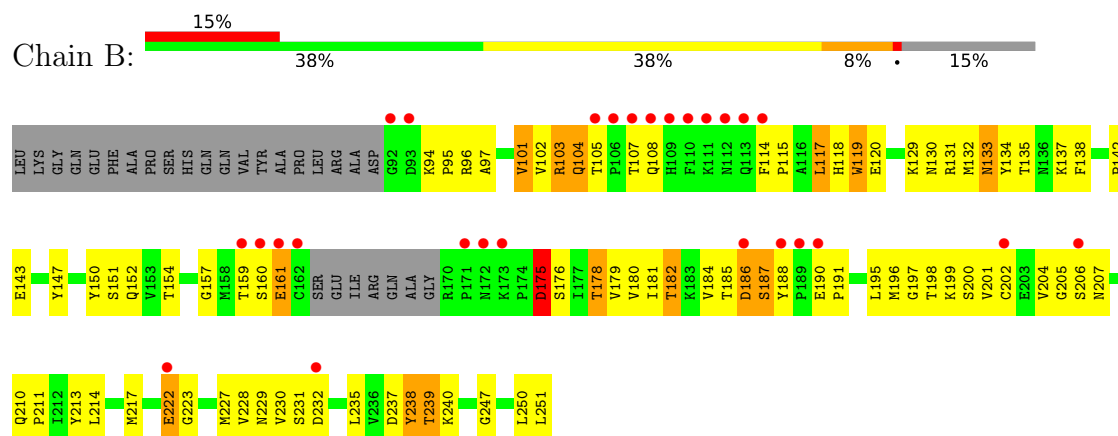
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

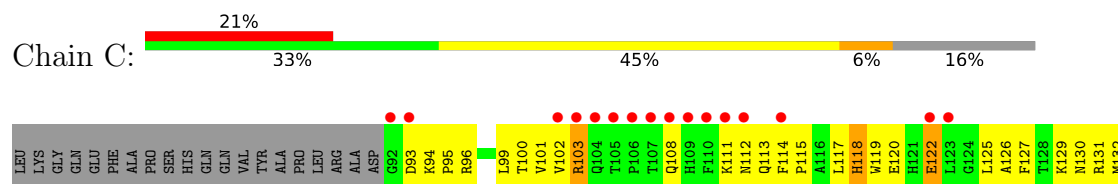
• Molecule 1: TNF superfamily ligand TL1A

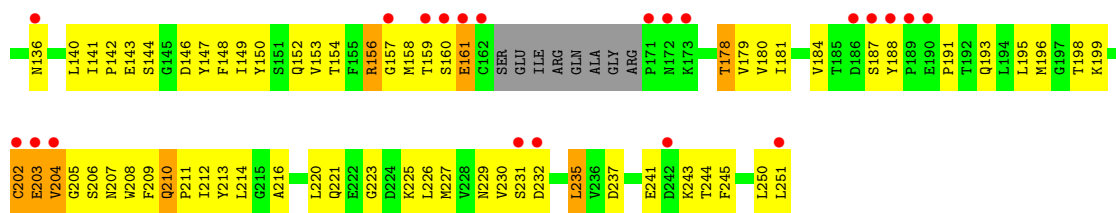


• Molecule 1: TNF superfamily ligand TL1A



• Molecule 1: TNF superfamily ligand TL1A





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	116.75Å 116.75Å 118.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 3.00 40.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	89.2 (40.00-3.00) 89.2 (40.00-3.00)	Depositor EDS
R_{merge}	0.31	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 3.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.243 , 0.295 0.244 , 0.295	Depositor DCC
R_{free} test set	828 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 79.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.033 for -h,l,k 0.023 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	3591	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.49	0/1232	0.95	2/1675 (0.1%)
1	B	0.50	0/1212	0.98	6/1648 (0.4%)
1	C	0.47	0/1201	0.89	0/1637
All	All	0.49	0/3645	0.94	8/4960 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	119	TRP	N-CA-C	9.19	120.37	108.34
1	A	239	THR	N-CA-C	6.36	119.03	111.33
1	B	119	TRP	N-CA-C	5.95	117.62	109.18
1	B	186	ASP	N-CA-C	-5.59	105.09	111.07
1	B	103	ARG	N-CA-C	5.49	117.27	111.28
1	B	120	GLU	N-CA-C	-5.48	100.31	109.24
1	B	175	ASP	N-CA-C	-5.43	102.42	110.46
1	B	133	ASN	N-CA-C	5.00	116.84	107.99

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1200	0	1149	79	0
1	B	1183	0	1127	78	0
1	C	1170	0	1100	95	0
2	A	1	0	0	0	0
3	A	16	0	0	0	0
3	B	12	0	0	1	0
3	C	9	0	0	2	0
All	All	3591	0	3376	235	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 34.

All (235) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:178:THR:HB	1:C:198:THR:HG22	1.24	1.08
1:A:185:THR:HG22	1:A:187:SER:H	1.29	0.95
1:C:178:THR:HG23	1:C:231:SER:HB3	1.47	0.93
1:A:183:LYS:HB2	1:A:194:LEU:HD11	1.51	0.92
1:C:232:ASP:HB3	1:C:235:LEU:HD22	1.52	0.89
1:C:181:ILE:HD12	1:C:195:LEU:HD12	1.58	0.85
1:B:117:LEU:HD11	1:B:230:VAL:HG13	1.59	0.84
1:A:105:THR:H	1:A:118:HIS:HE1	1.26	0.84
1:C:178:THR:CB	1:C:198:THR:HG22	2.08	0.83
1:B:118:HIS:HB3	1:B:137:LYS:HE3	1.62	0.82
1:B:105:THR:H	1:B:118:HIS:HE1	1.24	0.81
1:C:117:LEU:HD11	1:C:230:VAL:HG13	1.60	0.81
1:C:227:MET:HE3	1:C:229:ASN:HD21	1.46	0.81
1:B:178:THR:HG23	1:B:231:SER:HB3	1.63	0.79
1:B:94:LYS:HB3	1:B:250:LEU:HB3	1.64	0.78
1:A:178:THR:HG22	1:A:231:SER:HB3	1.67	0.77
1:B:190:GLU:OE1	1:B:191:PRO:HD2	1.86	0.76
1:A:143:GLU:HA	1:A:222:GLU:HG2	1.67	0.76
1:C:94:LYS:HB2	1:C:95:PRO:HD2	1.67	0.76
1:A:200:SER:HB3	1:C:209:PHE:CE2	2.22	0.75
1:C:178:THR:HB	1:C:198:THR:CG2	2.13	0.75
1:C:193:GLN:HE22	1:C:196:MET:HE3	1.53	0.74
1:B:204:VAL:HG23	1:C:202:CYS:SG	2.27	0.74
1:A:243:LYS:HE3	1:B:196:MET:HE3	1.69	0.73
1:C:195:LEU:HD21	1:C:216:ALA:HB3	1.70	0.72
1:A:105:THR:H	1:A:118:HIS:CE1	2.12	0.67
1:A:114:PHE:HB3	1:A:229:ASN:HB3	1.74	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ILE:HG21	1:A:181:ILE:HG21	1.76	0.67
1:B:105:THR:H	1:B:118:HIS:CE1	2.10	0.67
1:C:195:LEU:HB3	1:C:214:LEU:HD13	1.76	0.67
1:B:181:ILE:HD12	1:B:195:LEU:HD12	1.77	0.66
1:A:104:GLN:HE22	1:A:118:HIS:H	1.43	0.66
1:B:133:ASN:C	1:B:133:ASN:HD22	2.03	0.65
1:B:103:ARG:HA	1:B:238:TYR:CD1	2.31	0.65
1:C:95:PRO:HB2	1:C:132:MET:HE3	1.78	0.65
1:C:147:TYR:CE1	1:C:250:LEU:HD13	2.32	0.64
1:B:143:GLU:HA	1:B:222:GLU:HB2	1.79	0.64
1:C:159:THR:H	1:C:207:ASN:HA	1.62	0.64
1:A:143:GLU:HA	1:A:222:GLU:CG	2.28	0.64
1:A:159:THR:HG23	1:A:203:GLU:O	1.98	0.63
1:B:115:PRO:HD2	1:B:230:VAL:O	1.98	0.63
1:B:103:ARG:HA	1:B:238:TYR:HD1	1.63	0.63
1:B:103:ARG:O	1:B:104:GLN:HB2	1.99	0.61
1:C:193:GLN:NE2	1:C:196:MET:HE3	2.15	0.61
1:C:203:GLU:HG3	1:C:208:TRP:CD1	2.35	0.61
1:C:195:LEU:HD21	1:C:216:ALA:CB	2.30	0.61
1:C:120:GLU:OE2	1:C:122:GLU:HG2	2.01	0.60
1:B:117:LEU:N	1:B:117:LEU:HD12	2.16	0.59
1:A:155:PHE:CG	1:A:177:ILE:HD12	2.38	0.59
1:B:129:LYS:HG3	1:B:130:ASN:ND2	2.17	0.59
1:A:108:GLN:HE21	1:A:109:HIS:N	2.01	0.59
1:A:108:GLN:HE21	1:A:108:GLN:C	2.11	0.59
1:B:199:LYS:HD3	3:B:252:HOH:O	2.01	0.58
1:B:95:PRO:HA	1:B:130:ASN:HB2	1.84	0.58
1:A:175:ASP:HB2	1:A:201:VAL:HG13	1.85	0.58
1:A:127:PHE:CZ	1:B:217:MET:HE3	2.39	0.58
1:A:143:GLU:CA	1:A:222:GLU:HG2	2.32	0.58
1:C:157:GLY:O	1:C:207:ASN:HB2	2.04	0.57
1:B:96:ARG:HH21	1:C:146:ASP:HB2	1.69	0.57
1:A:131:ARG:O	1:A:142:PRO:HB3	2.04	0.57
1:B:133:ASN:HD22	1:B:134:TYR:N	2.02	0.57
1:B:157:GLY:O	1:B:207:ASN:HB2	2.04	0.57
1:C:100:THR:CG2	1:C:241:GLU:HB2	2.35	0.57
1:B:114:PHE:CE2	1:B:231:SER:HA	2.40	0.57
1:C:156:ARG:HD2	1:C:237:ASP:HB2	1.86	0.56
1:C:144:SER:OG	1:C:221:GLN:HA	2.05	0.56
1:A:148:PHE:C	1:A:149:ILE:HD13	2.29	0.56
1:C:112:ASN:CG	1:C:113:GLN:H	2.14	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ILE:HD13	1:A:149:ILE:N	2.20	0.56
1:B:239:THR:HG22	1:B:240:LYS:N	2.19	0.56
1:C:220:LEU:HD13	1:C:226:LEU:HD11	1.88	0.56
1:B:180:VAL:HG12	1:B:182:THR:HG22	1.88	0.56
1:A:183:LYS:HB2	1:A:194:LEU:CD1	2.32	0.56
1:C:193:GLN:HE22	1:C:196:MET:CE	2.18	0.55
1:B:117:LEU:HD12	1:B:117:LEU:H	1.70	0.55
1:C:203:GLU:HG3	1:C:208:TRP:HB3	1.88	0.55
1:A:104:GLN:HA	1:A:118:HIS:CE1	2.42	0.55
1:A:240:LYS:HB2	1:A:243:LYS:HG3	1.89	0.55
1:A:95:PRO:HD2	1:A:250:LEU:HD22	1.89	0.55
1:B:103:ARG:HG3	1:B:238:TYR:CE1	2.41	0.55
1:C:96:ARG:HD2	1:C:129:LYS:HD2	1.89	0.55
1:C:99:LEU:HB3	1:C:119:TRP:HB3	1.88	0.55
1:A:176:SER:HB3	1:C:209:PHE:HZ	1.72	0.55
1:C:147:TYR:CZ	1:C:250:LEU:HD13	2.42	0.55
1:A:115:PRO:HD2	1:A:230:VAL:O	2.07	0.54
1:B:151:SER:HB3	1:B:181:ILE:HD11	1.89	0.54
1:C:141:ILE:CD1	1:C:226:LEU:HD13	2.37	0.54
1:C:180:VAL:HG22	1:C:196:MET:HG3	1.88	0.54
1:B:199:LYS:HD2	1:B:210:GLN:OE1	2.06	0.54
1:B:238:TYR:N	1:B:238:TYR:CD2	2.74	0.54
1:C:232:ASP:CB	1:C:235:LEU:HD22	2.32	0.54
1:B:101:VAL:HG22	1:B:102:VAL:H	1.73	0.54
1:B:96:ARG:HD2	1:C:146:ASP:OD1	2.08	0.54
1:A:197:GLY:N	1:A:214:LEU:HD11	2.23	0.53
1:B:160:SER:O	1:B:161:GLU:C	2.51	0.53
1:B:237:ASP:OD1	1:B:239:THR:HB	2.08	0.53
1:C:199:LYS:HD2	1:C:210:GLN:OE1	2.08	0.53
1:C:136:ASN:HB3	3:C:259:HOH:O	2.07	0.53
1:A:101:VAL:HB	1:A:244:THR:HB	1.90	0.53
1:A:185:THR:HG22	1:A:187:SER:N	2.12	0.53
1:B:118:HIS:HB3	1:B:137:LYS:CE	2.37	0.53
1:A:100:THR:O	1:A:119:TRP:HB3	2.09	0.52
1:C:118:HIS:HB3	3:C:254:HOH:O	2.09	0.52
1:B:131:ARG:O	1:B:142:PRO:HB3	2.10	0.52
1:C:117:LEU:HD12	1:C:117:LEU:N	2.25	0.52
1:B:197:GLY:HA3	1:B:214:LEU:HD21	1.92	0.52
1:C:100:THR:HB	1:C:241:GLU:HB2	1.91	0.51
1:A:173:LYS:HA	1:C:206:SER:HB3	1.91	0.51
1:C:209:PHE:O	1:C:210:GLN:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:PHE:CE2	1:B:200:SER:HB3	2.46	0.51
1:C:203:GLU:HG3	1:C:208:TRP:CG	2.46	0.51
1:C:101:VAL:HG22	1:C:102:VAL:N	2.26	0.51
1:C:140:LEU:HD11	1:C:223:GLY:HA2	1.93	0.51
1:B:117:LEU:HD11	1:B:230:VAL:CG1	2.37	0.50
1:B:186:ASP:O	1:B:187:SER:O	2.30	0.50
1:B:107:THR:O	1:B:107:THR:HG23	2.11	0.50
1:C:146:ASP:O	1:C:250:LEU:HD12	2.12	0.50
1:C:149:ILE:HD12	1:C:149:ILE:N	2.26	0.50
1:A:174:PRO:HB3	1:A:202:CYS:SG	2.52	0.50
1:B:96:ARG:NH2	1:C:146:ASP:HB2	2.26	0.50
1:C:184:VAL:HG22	1:C:191:PRO:HB3	1.92	0.50
1:A:104:GLN:NE2	1:A:117:LEU:HA	2.27	0.50
1:C:148:PHE:HB2	1:C:251:LEU:HD11	1.93	0.50
1:B:152:GLN:HB2	1:B:213:TYR:HD1	1.76	0.50
1:A:225:LYS:C	1:A:226:LEU:HD22	2.37	0.49
1:C:149:ILE:N	1:C:149:ILE:CD1	2.75	0.49
1:B:133:ASN:C	1:B:133:ASN:ND2	2.67	0.49
1:C:149:ILE:HG21	1:C:181:ILE:HG21	1.94	0.49
1:A:103:ARG:O	1:A:104:GLN:CB	2.60	0.49
1:A:154:THR:HA	1:A:211:PRO:HA	1.95	0.49
1:B:187:SER:OG	1:B:188:TYR:N	2.46	0.49
1:C:103:ARG:HE	1:C:103:ARG:HA	1.78	0.49
1:C:204:VAL:HG22	1:C:204:VAL:O	2.13	0.48
1:B:105:THR:N	1:B:118:HIS:HE1	2.01	0.48
1:C:95:PRO:HA	1:C:130:ASN:HB2	1.95	0.48
1:C:251:LEU:HD12	1:C:251:LEU:N	2.28	0.48
1:A:105:THR:N	1:A:118:HIS:HE1	2.03	0.48
1:C:125:LEU:HD12	1:C:241:GLU:HG3	1.96	0.48
1:A:138:PHE:HE1	1:A:227:MET:HE2	1.79	0.48
1:A:127:PHE:HZ	1:B:217:MET:HE3	1.78	0.48
1:B:232:ASP:CG	1:B:235:LEU:HD13	2.39	0.47
1:A:220:LEU:HD13	1:A:226:LEU:HD21	1.96	0.47
1:A:188:TYR:CD1	1:A:188:TYR:N	2.83	0.47
1:A:243:LYS:HE2	1:B:198:THR:HG21	1.96	0.47
1:C:112:ASN:CG	1:C:113:GLN:N	2.73	0.47
1:C:142:PRO:HG2	1:C:143:GLU:H	1.79	0.47
1:C:100:THR:HG21	1:C:241:GLU:HB2	1.96	0.47
1:A:131:ARG:HH22	1:A:143:GLU:HB3	1.80	0.47
1:A:204:VAL:O	1:A:204:VAL:HG23	2.14	0.47
1:C:141:ILE:HD11	1:C:226:LEU:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:ARG:HH21	1:C:156:ARG:HG3	1.80	0.46
1:A:187:SER:HB2	1:A:188:TYR:CD1	2.51	0.46
1:B:250:LEU:HD13	1:B:251:LEU:N	2.30	0.46
1:C:251:LEU:HD12	1:C:251:LEU:H	1.80	0.46
1:A:152:GLN:O	1:A:244:THR:HA	2.16	0.46
1:C:152:GLN:HB3	1:C:245:PHE:CZ	2.51	0.46
1:A:171:PRO:HG2	1:C:206:SER:HA	1.98	0.45
1:C:100:THR:CB	1:C:241:GLU:HB2	2.46	0.45
1:C:114:PHE:N	1:C:115:PRO:HD3	2.30	0.45
1:A:179:VAL:HA	1:A:229:ASN:O	2.17	0.45
1:A:211:PRO:HD3	1:B:198:THR:O	2.17	0.45
1:C:187:SER:OG	1:C:188:TYR:HD1	1.99	0.45
1:B:238:TYR:HD2	1:B:239:THR:N	2.15	0.45
1:C:95:PRO:HG2	1:C:250:LEU:HB2	1.98	0.45
1:A:233:ILE:HG22	1:A:234:SER:N	2.30	0.44
1:A:242:ASP:OD2	1:A:242:ASP:N	2.46	0.44
1:B:130:ASN:O	1:B:131:ARG:HB2	2.17	0.44
1:B:159:THR:HG22	1:B:161:GLU:HB3	1.99	0.44
1:C:153:VAL:HA	1:C:244:THR:HG23	2.00	0.44
1:A:119:TRP:CD1	1:A:119:TRP:N	2.85	0.44
1:A:155:PHE:CD1	1:A:177:ILE:HD12	2.52	0.44
1:A:208:TRP:O	1:A:209:PHE:HB3	2.18	0.44
1:B:150:TYR:CZ	1:B:247:GLY:HA3	2.52	0.44
1:B:180:VAL:HG13	1:B:195:LEU:O	2.17	0.44
1:C:154:THR:HA	1:C:211:PRO:HA	2.00	0.44
1:A:104:GLN:NE2	1:A:118:HIS:H	2.13	0.43
1:B:104:GLN:HA	1:B:118:HIS:CE1	2.53	0.43
1:C:126:ALA:O	1:C:127:PHE:HB3	2.18	0.43
1:A:104:GLN:HE22	1:A:118:HIS:N	2.13	0.43
1:A:104:GLN:HE22	1:A:117:LEU:HA	1.82	0.43
1:A:105:THR:HG23	1:A:106:PRO:HD2	1.99	0.43
1:B:97:ALA:HB2	1:B:132:MET:HG2	2.00	0.43
1:A:195:LEU:HD11	1:A:216:ALA:HB3	2.01	0.43
1:B:101:VAL:HG22	1:B:102:VAL:N	2.33	0.43
1:C:150:TYR:HA	1:C:214:LEU:O	2.18	0.43
1:B:181:ILE:HG12	1:B:228:VAL:HG22	2.00	0.43
1:C:204:VAL:O	1:C:204:VAL:HG13	2.17	0.43
1:A:95:PRO:HB3	1:A:142:PRO:HG3	2.00	0.43
1:A:162:CYS:HB3	1:A:172:ASN:CB	2.49	0.43
1:A:103:ARG:NH2	1:A:103:ARG:HB2	2.34	0.43
1:B:119:TRP:CD1	1:B:119:TRP:N	2.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:PHE:CE2	1:C:231:SER:HA	2.54	0.43
1:C:160:SER:O	1:C:161:GLU:C	2.61	0.43
1:A:139:LEU:HD12	1:A:139:LEU:HA	1.91	0.43
1:A:238:TYR:CD1	1:A:238:TYR:C	2.96	0.43
1:A:195:LEU:HD13	1:A:218:PHE:HZ	1.84	0.42
1:A:152:GLN:HB3	1:A:245:PHE:CE1	2.54	0.42
1:B:138:PHE:HE1	1:B:227:MET:HE2	1.84	0.42
1:A:200:SER:HB3	1:C:209:PHE:CZ	2.55	0.42
1:B:95:PRO:CB	1:B:142:PRO:HG3	2.50	0.42
1:B:147:TYR:CZ	1:B:250:LEU:HD23	2.54	0.42
1:C:152:GLN:HA	1:C:212:ILE:O	2.19	0.42
1:A:123:LEU:O	1:A:126:ALA:HB3	2.19	0.42
1:B:222:GLU:HG3	1:B:223:GLY:N	2.35	0.42
1:C:152:GLN:HB2	1:C:213:TYR:HD1	1.85	0.42
1:B:190:GLU:OE1	1:B:190:GLU:HA	2.19	0.41
1:C:99:LEU:HB3	1:C:119:TRP:CB	2.49	0.41
1:C:154:THR:HG21	1:C:243:LYS:HE2	2.02	0.41
1:A:102:VAL:O	1:A:103:ARG:C	2.63	0.41
1:C:131:ARG:O	1:C:142:PRO:HB3	2.20	0.41
1:C:225:LYS:HE3	1:C:225:LYS:HB2	1.82	0.41
1:A:181:ILE:HD12	1:A:181:ILE:N	2.34	0.41
1:C:220:LEU:HD11	1:C:226:LEU:HD21	2.02	0.41
1:C:156:ARG:HD3	1:C:235:LEU:O	2.20	0.41
1:A:118:HIS:HB3	1:A:137:LYS:HG2	2.01	0.41
1:A:185:THR:HG22	1:A:186:ASP:N	2.35	0.41
1:A:217:MET:HE1	1:C:96:ARG:HB2	2.01	0.41
1:C:118:HIS:ND1	1:C:118:HIS:N	2.68	0.41
1:B:154:THR:HA	1:B:211:PRO:HA	2.03	0.41
1:B:175:ASP:OD2	1:B:175:ASP:N	2.51	0.41
1:B:205:GLY:O	1:B:206:SER:C	2.64	0.41
1:B:182:THR:HG23	1:B:229:ASN:HD22	1.86	0.41
1:B:238:TYR:N	1:B:238:TYR:HD2	2.18	0.41
1:A:104:GLN:HA	1:A:118:HIS:HE1	1.84	0.40
1:B:152:GLN:HE22	1:C:214:LEU:HD21	1.86	0.40
1:C:159:THR:HB	1:C:205:GLY:O	2.21	0.40
1:A:96:ARG:HE	1:A:96:ARG:HB3	1.67	0.40
1:B:114:PHE:HB3	1:B:229:ASN:HB3	2.02	0.40
1:B:190:GLU:HA	1:B:191:PRO:HD2	1.87	0.40
1:C:203:GLU:HG3	1:C:208:TRP:CB	2.50	0.40
1:B:103:ARG:O	1:B:104:GLN:CB	2.67	0.40
1:B:185:THR:OG1	1:B:186:ASP:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:LYS:CB	1:C:95:PRO:HD2	2.44	0.40
1:C:158:MET:HA	1:C:207:ASN:CB	2.51	0.40
1:A:130:ASN:N	1:A:130:ASN:HD22	2.18	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	149/180 (83%)	133 (89%)	10 (7%)	6 (4%)	2	14
1	B	149/180 (83%)	129 (87%)	15 (10%)	5 (3%)	3	17
1	C	148/180 (82%)	120 (81%)	20 (14%)	8 (5%)	1	9
All	All	446/540 (83%)	382 (86%)	45 (10%)	19 (4%)	2	12

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	160	SER
1	A	187	SER
1	A	191	PRO
1	B	161	GLU
1	B	187	SER
1	C	111	LYS
1	C	203	GLU
1	C	204	VAL
1	A	103	ARG
1	A	209	PHE
1	B	108	GLN
1	B	239	THR
1	C	93	ASP
1	C	108	GLN

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Mol	Chain	Res	Type
1	C	202	CYS
1	B	104	GLN
1	C	122	GLU
1	A	104	GLN
1	C	210	GLN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	130/160 (81%)	115 (88%)	15 (12%)	5	24
1	B	127/160 (79%)	114 (90%)	13 (10%)	7	29
1	C	124/160 (78%)	117 (94%)	7 (6%)	19	52
All	All	381/480 (79%)	346 (91%)	35 (9%)	8	33

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	LEU
1	A	103	ARG
1	A	108	GLN
1	A	117	LEU
1	A	119	TRP
1	A	120	GLU
1	A	128	THR
1	A	140	LEU
1	A	149	ILE
1	A	179	VAL
1	A	180	VAL
1	A	186	ASP
1	A	195	LEU
1	A	203	GLU
1	A	233	ILE
1	B	101	VAL
1	B	117	LEU

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Mol	Chain	Res	Type
1	B	135	THR
1	B	175	ASP
1	B	176	SER
1	B	178	THR
1	B	179	VAL
1	B	182	THR
1	B	184	VAL
1	B	201	VAL
1	B	202	CYS
1	B	222	GLU
1	B	238	TYR
1	C	103	ARG
1	C	118	HIS
1	C	156	ARG
1	C	161	GLU
1	C	178	THR
1	C	179	VAL
1	C	235	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	108	GLN
1	A	118	HIS
1	A	130	ASN
1	A	193	GLN
1	A	221	GLN
1	B	113	GLN
1	B	118	HIS
1	B	133	ASN
1	B	136	ASN
1	B	152	GLN
1	B	210	GLN
1	B	229	ASN
1	C	108	GLN
1	C	112	ASN
1	C	130	ASN
1	C	136	ASN
1	C	193	GLN
1	C	221	GLN
1	C	229	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	153/180 (85%)	0.88	22 (14%) 6 3	14, 39, 93, 113	0
1	B	153/180 (85%)	1.06	27 (17%) 4 3	16, 37, 100, 124	0
1	C	152/180 (84%)	1.42	37 (24%) 2 1	20, 56, 118, 153	0
All	All	458/540 (84%)	1.12	86 (18%) 3 2	14, 44, 98, 153	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	108	GLN	7.3
1	C	104	GLN	5.9
1	C	110	PHE	5.5
1	B	110	PHE	5.3
1	B	109	HIS	5.0
1	A	172	ASN	4.8
1	B	172	ASN	4.7
1	B	188	TYR	4.7
1	C	107	THR	4.5
1	B	162	CYS	4.5
1	B	112	ASN	4.4
1	C	111	LYS	4.4
1	C	203	GLU	4.3
1	A	188	TYR	4.3
1	C	162	CYS	4.2
1	B	107	THR	4.2
1	B	108	GLN	4.1
1	B	111	LYS	4.0
1	C	172	ASN	4.0
1	C	106	PRO	3.7
1	C	189	PRO	3.7
1	C	188	TYR	3.7
1	A	173	LYS	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	171	PRO	3.6
1	A	160	SER	3.6
1	A	186	ASP	3.6
1	A	189	PRO	3.6
1	C	92	GLY	3.6
1	B	189	PRO	3.6
1	B	159	THR	3.5
1	C	109	HIS	3.5
1	C	93	ASP	3.5
1	A	162	CYS	3.5
1	B	160	SER	3.3
1	C	159	THR	3.3
1	C	187	SER	3.2
1	C	105	THR	3.2
1	B	92	GLY	3.2
1	B	105	THR	3.2
1	B	206	SER	3.1
1	B	93	ASP	3.1
1	A	204	VAL	3.1
1	A	170	ARG	3.0
1	C	122	GLU	3.0
1	B	106	PRO	3.0
1	B	171	PRO	3.0
1	C	171	PRO	3.0
1	A	202	CYS	2.9
1	C	251	LEU	2.9
1	B	202	CYS	2.8
1	C	202	CYS	2.8
1	A	224	ASP	2.7
1	C	160	SER	2.7
1	A	175	ASP	2.7
1	C	112	ASN	2.7
1	B	113	GLN	2.7
1	A	232	ASP	2.7
1	B	114	PHE	2.6
1	A	190	GLU	2.6
1	A	93	ASP	2.5
1	A	161	GLU	2.5
1	C	186	ASP	2.5
1	C	102	VAL	2.5
1	C	103	ARG	2.4
1	B	161	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	191	PRO	2.4
1	A	185	THR	2.4
1	C	204	VAL	2.4
1	C	231	SER	2.3
1	C	232	ASP	2.3
1	C	190	GLU	2.3
1	B	190	GLU	2.2
1	C	242	ASP	2.2
1	C	136	ASN	2.2
1	A	146	ASP	2.2
1	B	173	LYS	2.2
1	C	123	LEU	2.1
1	A	222	GLU	2.1
1	B	232	ASP	2.1
1	B	222	GLU	2.1
1	C	114	PHE	2.1
1	B	186	ASP	2.1
1	C	173	LYS	2.1
1	A	108	GLN	2.1
1	C	157	GLY	2.1
1	C	161	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	A	271	1/1	0.96	0.18	6,6,6,6	0

6.5 Other polymers [i](#)

There are no such residues in this entry.